

M	Final Report
	Revised Report

Report Date: 24-May-17 11:45

Laboratory Report SC34648

Gulf Oil L.P. 281 Eastern Avenue Chelsea, MA 02150 Attn: Andrew P. Adams

Project: Gulf Terminal - Chelsea, MA

Project #: Gulf Chelsea

I attest that the information contained within the report has been reviewed for accuracy and checked against the quality control requirements for each method. These results relate only to the sample(s) as received.

All applicable NELAC requirements have been met.

Massachusetts # M-MA138/MA1110 Connecticut # PH-0777 Florida # E87936 Maine # MA138 New Hampshire # 2972/2538 New Jersey # MA011 New York # 11393 Pennsylvania # 68-04426/68-02924 Rhode Island # LAO00348 USDA # P330-15-00375 Vermont # VT-11393



Authorized by:

Christina White Laboratory Director

Christina a. White

Eurofins Spectrum Analytical holds primary certification in the State of Massachusetts for the analytes as indicated with an X in the "Cert." column within this report. Please note that the State of Massachusetts does not offer certification for all analytes. Please refer to our website for specific certification holdings in each state.

Please note that this report contains 12 pages of analytical data plus Chain of Custody document(s). When the Laboratory Report is indicated as revised, this report supersedes any previously dated reports for the laboratory ID(s) referenced above. Where this report identifies subcontracted analyses, copies of the subcontractor's test report are available upon request. This report may not be reproduced, except in full, without written approval from Eurofins Spectrum Analytical, Inc.

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Please contact the Laboratory or Technical Director at 800-789-9115 with any questions regarding the data contained in this laboratory report.

Sample Summary

Work Order: SC34648

Project: Gulf Terminal - Chelsea, MA

Project Number: Gulf Chelsea

Laboratory IDClient Sample IDMatrixDate SampledDate ReceivedSC34648-01Outfall 003Surface Water13-May-17 21:0015-May-17 14:05

CASE NARRATIVE:

Data has been reported to the RDL. This report excludes estimated concentrations detected below the RDL and above the MDL (J-Flag).

All non-detects and all results below the reporting limit are reported as "<" (less than) the reporting limit in this report.

The samples were received 3.8 degrees Celsius, please refer to the Chain of Custody for details specific to temperature upon receipt. An infrared thermometer with a tolerance of +/- 1.0 degrees Celsius was used immediately upon receipt of the samples.

If a Matrix Spike (MS), Matrix Spike Duplicate (MSD) or Duplicate (DUP) was not requested on the Chain of Custody, method criteria may have been fulfilled with a source sample not of this Sample Delivery Group.

Analyses for Total Hardness, pH, and Total Residual Chlorine fall under the state of Pennsylvania code Chapter 252.6 accreditation by rule.

See below for any non-conformances and issues relating to quality control samples and/or sample analysis/matrix.

SW846 8260C

Calibration:

1705009

Analyte quantified by quadratic equation type calibration.

Naphthalene

This affected the following samples:

1708138-BLK1

1708138-BLK2

1708138-BS1

1708138-BS2

1708138-BSD1

1708138-BSD2

Outfall 003

S704360-ICV1

S704555-CCV1

Laboratory Control Samples:

1708138-BS1

LCS/LCSD were analyzed in place of MS/MSD.

1708138-BSD1

LCS/LCSD were analyzed in place of MS/MSD.

SW846 8270D SIM

Calibration:

1704025

Analyte quantified by quadratic equation type calibration.

Benzo (a) pyrene

Benzo (e) pyrene-d12

This laboratory report is not valid without an authorized signature on the cover page.

SW846 8270D SIM

Calibration:

1704025

This affected the following samples:

1708203-BLK2

1708203-BS2

1708203-BSD2

Outfall 003

S703654-ICV1

S704678-CCV1

S704750-CCV1

Laboratory Control Samples:

1708203 BSD

Benzo (a) pyrene RPD 25% (20%) is outside individual acceptance criteria.

Naphthalene RPD 21% (20%) is outside individual acceptance criteria.

1708203-BSD2

RPD out of acceptance range. The batch is accepted based upon LCS and/or LCSD recovery.

Benzo (a) pyrene

Naphthalene

Phenanthrene

Pyrene

24-May-17 11:45 Page 4 of 12

Sample Acceptance Check Form

Project:	Gulf Terminal - Chelsea, MA / Gulf Chelsea			
Work Order:	SC34648			
Sample(s) received on:	5/15/2017			
The following outlines th	he condition of samples for the attached Chain of Custody upon receipt.			
		<u>Yes</u>	<u>No</u>	N/A
Were custody se	als present?		\checkmark	
Were custody se	als intact?			\checkmark
Were samples re	ceived at a temperature of $\leq 6^{\circ}$ C?	\checkmark		
Were samples re	frigerated upon transfer to laboratory representative?	\checkmark		
Were sample con	ntainers received intact?	\checkmark		
	operly labeled (labels affixed to sample containers and include sample ID, site project number and the collection date)?			
Were samples ac	companied by a Chain of Custody document?	\checkmark		
include sample I	Sustody document include proper, full, and complete documentation, which shall D, site location, and/or project number, date and time of collection, collector's name, e, sample matrix and any special remarks concerning the sample?	$ \checkmark $		
Did sample cont	ainer labels agree with Chain of Custody document?	\checkmark		
Were samples re	ceived within method-specific holding times?	\checkmark		

Client:

Gulf Oil L.P.

Summary of Hits

Lab ID: SC34648-01

Client ID: Outfall 003

Parameter	Result	Flag	Reporting Limit	Units	Analytical Method
Total Suspended Solids	9.8		1.0	mg/l	SM2540D (11)
Naphthalene	0.076		0.051	μg/l	SW846 8270D SIM

Please note that because there are no reporting limits associated with hazardous waste characterizations or micro analyses, this summary does not include hits from these analyses if included in this work order.

Sample Id Outfall 06 SC34648-					Project # Chelsea		<u>Matrix</u> Surface W		ection Date -May-17 21			eceived May-17	
CAS No.	Analyte(s)	Result	Flag	Units	*RDL	MDL	Dilution	Method Ref.	Prepared	Analyzed	Analyst	Batch	Cert.
Volatile O	rganic Compounds												
Volatile O	rganic Aromatics by SW8	46 8260											
	by method SW846 5030												
71-43-2	Benzene	< 1.0		μg/l	1.0	0.3	1	SW846 8260C	16-May-1 7	17-May-1 7	GMA	1708138	
91-20-3	Naphthalene	< 1.0		μg/l	1.0	0.4	1	"	"	"	"	"	
Surrogate	recoveries:												
460-00-4	4-Bromofluorobenzene	96			70-13	80 %		"	"	"	"	"	
2037-26-5	Toluene-d8	98			70-13	80 %		"	"	"	"	"	
17060-07-0	1,2-Dichloroethane-d4	102			70-13	80 %		n	"	"	"	"	
1868-53-7	Dibromofluoromethane	98			70-13	80 %		"	"	"	"	"	
Semivolati	ile Organic Compounds by	GCMS											
SVOCs by	y SIM by method SW846 35100												
50-32-8	Benzo (a) pyrene	< 0.051		μg/l	0.051	0.006	1	SW846 8270D SIM	17-May-1 7	22-May-1 7	MSL	1708203	
91-20-3	Naphthalene	0.076		μg/l	0.051	0.007	1	"	"	"	"	"	
Surrogate	recoveries:												
205440-82-0	Benzo (e) pyrene-d12	69			30-13	80 %		"	"	u	"	"	
	le Petroleum Hydrocarbon by method SW846 35100												
	Oil & Grease	< 1.00		mg/l	1.00	0.915	1	EPA 1664B	22-May-1 7	22-May-1 7	KK	1708456	Χ
General C	hemistry Parameters												
	рН	9.16		pH Units			1	ASTM D 1293-99B	15-May-1 7 18:00	15-May-1 7 18:20	BD	1708122	Х
	Total Suspended Solids	9.8		mg/l	1.0	0.4	1	SM2540D (11)	16-May-1 7	17-May-1 7	CMB	1708147	Х

24-May-17 11:45 Page 7 of 12

Volatile Organic Compounds - Quality Control

	D L	El	TT 14	*DDI	Spike	Source	0/DEC	%REC	DDD	RPI
nalyte(s)	Result	Flag	Units	*RDL	Level	Result	%REC	Limits	RPD	Lim
W846 8260C										
atch 1708138 - SW846 5030 Water MS										
Blank (1708138-BLK1)					Pre	epared & Ar	nalyzed: 16	-May-17		
Benzene	< 1.0		μg/l	1.0						
Naphthalene	< 1.0		μg/l	1.0						
Surrogate: 4-Bromofluorobenzene	47.2		μg/l		50.0		94	70-130		
Surrogate: Toluene-d8	48.1		μg/l		50.0		96	70-130		
Surrogate: 1,2-Dichloroethane-d4	50.8		μg/l		50.0		102	70-130		
Surrogate: Dibromofluoromethane	48.0		μg/l		50.0		96	70-130		
Blank (1708138-BLK2)					Pre	epared & Ar	nalyzed: 16	-May-17		
Benzene	< 5.0	D	μg/l	5.0						
Naphthalene	< 5.0	D	μg/l	5.0						
Surrogate: 4-Bromofluorobenzene	46.5		μg/l		50.0		93	70-130		
Surrogate: Toluene-d8	48.1		μg/l		50.0		96	70-130		
Surrogate: 1,2-Dichloroethane-d4	51.5		μg/l		50.0		103	70-130		
Surrogate: Dibromofluoromethane	48.4		μg/l		50.0		97	70-130		
LCS (1708138-BS1)		QM10			Pre	epared & Ar	nalyzed: 16	-May-17		
Benzene	19.3		μg/l		20.0		97	70-130		
Naphthalene	22.1		μg/l		20.0		110	70-130		
Surrogate: 4-Bromofluorobenzene	49.0		μg/l		50.0		98	70-130		
Surrogate: Toluene-d8	48.7		μg/l		50.0		97	70-130		
Surrogate: 1,2-Dichloroethane-d4	49.6		μg/l		50.0		99	70-130		
Surrogate: Dibromofluoromethane	48.0		μg/l		50.0		96	70-130		
LCS (1708138-BS2)					Pre	epared: 16-	May-17 Ar	nalyzed: 17-N	<u> 1ay-17</u>	
Benzene	20.6	D	μg/l		20.0	•	103	70-130	 _	
Naphthalene	24.1	D	μg/l		20.0		121	70-130		
Surrogate: 4-Bromofluorobenzene	48.6		μg/l		50.0		97	70-130		
Surrogate: Toluene-d8	49.2		μg/l		50.0		98	70-130		
Surrogate: 1,2-Dichloroethane-d4	49.5		μg/l		50.0		99	70-130		
Surrogate: Dibromofluoromethane	48.3		μg/l		50.0		97	70-130		
LCS Dup (1708138-BSD1)		QM10			Pre	epared & Ar	nalyzed: 16	-May-17		
Benzene	21.1		μg/l		20.0		105	70-130	9	20
Naphthalene	23.4		μg/l		20.0		117	70-130	6	20
Surrogate: 4-Bromofluorobenzene	48.8		μg/l		50.0		98	70-130		
Surrogate: Toluene-d8	48.8		μg/l		50.0		98	70-130		
Surrogate: 1,2-Dichloroethane-d4	49.8		μg/l		50.0		100	70-130		
Surrogate: Dibromofluoromethane	47.5		μg/l		50.0		95	70-130		
LCS Dup (1708138-BSD2)					Pre	epared: 16-	May-17 Ar	nalyzed: 17-N	1ay-17	
Benzene	19.8	D	μg/l		20.0		99	70-130	4	20
Naphthalene	23.9	D	μg/l		20.0		119	70-130	0.9	20
Surrogate: 4-Bromofluorobenzene	49.0		μg/l		50.0		98	70-130		
Surrogate: Toluene-d8	49.1		μg/l		50.0		98	70-130		
Surrogate: 1,2-Dichloroethane-d4	48.9		μg/l		50.0		98	70-130		
Surrogate: Dibromofluoromethane	48.0		μg/l		50.0		96	70-130		

Semivolatile Organic Compounds by GCMS - Quality Control

1.7	D 1:	T.I.	**	*DDI	Spike	Source	0/DEC	%REC	n nn	RPD
nalyte(s)	Result	Flag	Units	*RDL	Level	Result	%REC	Limits	RPD	Limi
W846 8270D SIM										
atch 1708203 - SW846 3510C										
Blank (1708203-BLK2)					Pr	epared: 17-	May-17 Ar	nalyzed: 18-N	May-17	
Benzo (a) pyrene	< 0.050		μg/l	0.050						
Naphthalene	< 0.050		μg/l	0.050						
Phenanthrene	< 0.050		μg/l	0.050						
Pyrene	< 0.050		μg/l	0.050						
Surrogate: Benzo (e) pyrene-d12	0.780		μg/l		1.00		78	30-130		
LCS (1708203-BS2)					Pr	epared: 17-	May-17 Ar	nalyzed: 18-N	May-17	
Benzo (a) pyrene	0.721		μg/l	0.050	1.00		72	40-140		
Naphthalene	0.741		μg/l	0.050	1.00		74	40-140		
Phenanthrene	0.643		μg/l	0.050	1.00		64	40-140		
Pyrene	0.749		μg/l	0.050	1.00		75	40-140		
Surrogate: Benzo (e) pyrene-d12	0.800		μg/l		1.00		80	30-130		
LCS Dup (1708203-BSD2)					<u>Pr</u>	epared: 17-	May-17 Ar	nalyzed: 18-N	<u> May-17</u>	
Benzo (a) pyrene	0.929	QR9	μg/l	0.050	1.00		93	40-140	25	20
Naphthalene	0.916	QR9	μg/l	0.050	1.00		92	40-140	21	20
Phenanthrene	0.826	QR9	μg/l	0.050	1.00		83	40-140	25	20
Pyrene	0.977	QR9	μg/l	0.050	1.00		98	40-140	26	20
Surrogate: Benzo (e) pyrene-d12	0.990		μg/l		1.00		99	30-130		

Extractable Petroleum Hydrocarbons - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
<u>EPA 1664B</u>										
Batch 1708456 - SW846 3510C										
Blank (1708456-BLK1)					Pre	epared & An	nalyzed: 22-	May-17		
Oil & Grease	< 1.00		mg/l	1.00						
LCS (1708456-BS1)					Pre	epared & An	nalyzed: 22-	May-17		
Oil & Grease	13.4		mg/l	1.00	16.5		81	78-114		

General Chemistry Parameters - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
ASTM D 1293-99B										
Batch 1708122 - General Preparation										
<u>Duplicate (1708122-DUP1)</u>			Source: SC	<u>34648-01</u>	Pre	epared & An	alyzed: 15	5-May-17		
pH	9.16		pH Units			9.16			0	5
Reference (1708122-SRM1)					Pre	epared & An	alyzed: 15	5-May-17		
рH	6.01		pH Units		6.00		100	97.5-102. 5		
Reference (1708122-SRM2)					Pre	epared & An	alyzed: 15	5-May-17		
pН	6.02		pH Units		6.00		100	97.5-102. 5		
SM2540D (11)										
Batch 1708147 - General Preparation										
Blank (1708147-BLK1)					Pre	epared: 16-N	May-17 A	nalyzed: 17-M	lay-17	
Total Suspended Solids	< 0.5		mg/l	0.5						
LCS (1708147-BS1)					Pre	epared: 16-N	May-17 A	nalyzed: 17-M	lay-17	
Total Suspended Solids	98.0		mg/l	10.0	100		98	90-110		

Notes and Definitions

D Data reported from a dilution

QM10 LCS/LCSD were analyzed in place of MS/MSD.

QR9 RPD out of acceptance range. The batch is accepted based upon LCS and/or LCSD recovery.

dry Sample results reported on a dry weight basis

NR Not Reported

RPD Relative Percent Difference

OG The required Matrix Spike and Matrix Spike Duplicate (MS/MSD) for Oil & Grease method 1664B can only be analyzed

when the client has submitted sufficient sample volume. An extra liter per MS/MSD is required to fulfill the method QC criteria. Please refer to Chain of Custody and QC Summary (MS/MSD) of the Laboratory Report to verify ample sample

volume was submitted to fulfill the requirement.

pH The method for pH does not stipulate a specific holding time other than to state that the samples should be analyzed as

soon as possible. For aqueous samples the 40 CFR 136 specifies a holding time of 15 minutes from sampling to analysis. Therefore all aqueous pH samples not analyzed in the field are considered out of hold time at the time of sample receipt.

All soil samples are analyzed as soon as possible after sample receipt.

<u>Laboratory Control Sample (LCS)</u>: A known matrix spiked with compound(s) representative of the target analytes, which is used to document laboratory performance.

Matrix Duplicate: An intra-laboratory split sample which is used to document the precision of a method in a given sample matrix.

<u>Matrix Spike</u>: An aliquot of a sample spiked with a known concentration of target analyte(s). The spiking occurs prior to sample preparation and analysis. A matrix spike is used to document the bias of a method in a given sample matrix.

<u>Method Blank</u>: An analyte-free matrix to which all reagents are added in the same volumes or proportions as used in sample processing. The method blank should be carried through the complete sample preparation and analytical procedure. The method blank is used to document contamination resulting from the analytical process.

<u>Method Detection Limit (MDL)</u>: The minimum concentration of a substance that can be measured and reported with 99% confidence that the analyte concentration is greater than zero and is determined from analysis of a sample in a given matrix type containing the analyte.

Reportable Detection Limit (RDL): The lowest concentration that can be reliably achieved within specified limits of precision and accuracy during routine laboratory operating conditions. For many analytes the RDL analyte concentration is selected as the lowest non-zero standard in the calibration curve. While the RDL is approximately 5 to 10 times the MDL, the RDL for each sample takes into account the sample volume/weight, extract/digestate volume, cleanup procedures and, if applicable, dry weight correction. Sample RDLs are highly matrix-dependent.

<u>Surrogate</u>: An organic compound which is similar to the target analyte(s) in chemical composition and behavior in the analytical process, but which is not normally found in environmental samples. These compounds are spiked into all blanks, standards, and samples prior to analysis. Percent recoveries are calculated for each surrogate.

<u>Continuing Calibration Verification:</u> The calibration relationship established during the initial calibration must be verified at periodic intervals. Concentrations, intervals, and criteria are method specific.

CHAIN OF CUSTODY RECORD

□ Rush TAT - Date Needed:

☑ Standard TAT - 7 to 10 business days

Special Handling:

617 884 5980	Chalcas MA 02150 Wellesley, MA 02481-3705 Location: 281 Easte	281 Eastern Ave 80 William St, Suite 400	Gulf Oil LP Site Name:	Report To: Andrew Adams Invoice To: Christopher Gill Project No:	HANDAW I BULHANDUNI	1	SPECTRUM ANALYTICAL, INC. Page 1 of 1 Min. 24-hr n
Sampler(s):	Location:		Site Name:	Project No:			
	281 Eastern Ave, Chelsea State:		Gulf Chelsea Terminal	Gulf Chelsea		Samples disposed after 60 days unless otherwise instruc-	All TATs subject to laboratory approval Min. 24-hr notification needed for rushes

Featuring HANIBAL TECHNOLOGY		Se	Samples disposed after 60 days unless otherwise instructed.
Report To: Andrew Adams	Invoice To: Christopher Gill	Project No:	Gulf Chelsea
Gulf Oil LP	Gulf Oil LP		Gulf Chelses Terminal
281 Eastern Ave	80 William St, Suite 400	Site Name.	Cuit Original Formillar
Chelsea, MA 02150	Wellesley, MA 02481-3705	Location	281 Eastern Ave, Chelsea State: MA
Telephone #: 617.884.5980		Sampler(s):	
Project Mgr: Andrew Adams	P.O No.: Quote/RQN:		Anore Agams
F=Field Filtered 1=Na ₂ S2O ₃ 2=HCl 3=H ₂ SO ₄ 4=HNO ₃ 5=NaOH 6=Ascorbic Acid	4=HNO ₃ 5=NaOH 6=Ascorbic Acid	List Preservative Code below:	OA/OC Reporting Notes:
7=CH3OH 8 =NaHSO ₄ 9 =Deionized Water 10 =H ₃ PO ₄	11= none 12=		

* additional charges may appoply		none 12=	H3OH 8=NaHSO, 9=Deionized Water 10=H ₃ PO ₄ 11=
QA/QC Reporting Notes:	List Preservative Code below:	S	ield Filtered 1=Na ₂ S2O ₃ 2=HCl 3=H ₂ SO ₄ 4=HNO ₃ 5=NaOH 6=Ascorbic Acid
		F.O INC.	ect Mgr. Andrew Adams
DARE MARY	Andre	DO NO.	
	Sampler(s):		617 884 5980
281 Eastern Ave, Chelsea State: MA	Location: 281 Eastern	Wellesley, MA 02481-3705	Chelsea, MA 02150
186		80 William St, Suite 400	281 Eastern Ave
Gulf Chelsea Terminal	Site Name:	Gulf Oil LP	Gulf Oil LP
Gull Clieisea	Project No:	Invoice To: Christopher Gill	ort To: Andrew Adams

8=NaH	red	-	1		helsea, l
8=NaHSO ₄ 9=Deionized Water 10=H ₃ PO.	1=Na ₂ S2O ₃				helsea, MA 02150
onized Water	2=HC1		Andrew Adams	617.884.5980	
· 10=H ₃ PO ₄	$3=H_2SO_4$		dams	980	
	4=HNO ₃				
11= non	5=NaOH		P.		
11= none 12=	red 1=Na ₂ S ₂ O ₃ , 2 =HCl 3 =H ₂ SO ₄ 4 =HNO ₃ 5 =NaOH 6 =Ascorbic Acid		P.O No.:		Wellesley, MA 02481-3705
			Quote/RQN:		181-3705
	List Preservative Code below:		And	Sampler(s):	Location: 281 Ea
* additional charges may appply	QA/QC Reporting Notes:		STANON WAS		astern Ave, Chelsea State: MA

Project Mgr: Andrew Adams		P.O No.:	Quote/RQN:	And	Thorew Mound
F=Field Filtered 1=Na ₂ S2O ₃ 2=HCl 3	4=HNO ₃	6=Ascorbic Acid	, ,	List Preservative Code below:	QA/QC Reporting Notes:
/-CHECKLE OF MALESCA 2 Detonized frame to take 4				11 3 2 11	. annumum cumges may upppy
DW =Dinking Water GW =Groundwater S	SW=Surface Water WW=Waste Water	Water	Containers	Analysis	MA DEP MCP CAM Report?
O=Oil SO=Soil SL=Sludge A=Indoo	A=Indoor/Ambient Air SG=Soil Gas	,	8.		St.
X1=X2=_	X3=		Glass	e) (a) pyr	
G= Grab	C=Compsite	/pe itrix	Ambe Clear Plastic		
Lab ID: Sample ID:	Date: Time:	Ty	# of	napl PAI	State-specific reporting standards:
34642 1 Outfall 003	5-13-17 2166	6 G SW		×	
Outfall 003	5-13-17 2100	G SW		×	Required MLs:
Outfall 003	5-13-17 2100	G SW	ш	×	□ benzène 2 μg/L
Outfall 003	5-13-17 2100	G SW		×	naphthalene 5 µg/L
					□ benzo(a)pyrene 0.1 µg/L

Received by:

Date:

Time:

Temp °C

EDD format: E-mail to:

S

aadams@gulfoil.com, cgill@gulfoil.com

Condition upon receipt Custody Seals:

☐ Ptesent

☐ Intact ☐ Broken

☐ Ambient ☐ Iced

Rufrigerated | DI VOA Frozen

☐ Soil Jar Frozen

Batch Summary

1708122

General Chemistry Parameters

1708122-DUP1 1708122-SRM1 1708122-SRM2

SC34648-01 (Outfall 003)

1708138

Volatile Organic Compounds

1708138-BLK1 1708138-BLK2 1708138-BS1 1708138-BS2 1708138-BSD1 1708138-BSD2

SC34648-01 (Outfall 003)

1708147

General Chemistry Parameters

1708147-BLK1 1708147-BS1

SC34648-01 (Outfall 003)

1708203

Semivolatile Organic Compounds by GCMS

1708203-BLK2 1708203-BS2 1708203-BSD2

SC34648-01 (Outfall 003)

1708456

Extractable Petroleum Hydrocarbons

1708456-BLK1 1708456-BS1

SC34648-01 (Outfall 003)

S703654

Semivolatile Organic Compounds by GCMS

S703654-CAL1

S703654-CAL2

S703654-CAL3

S703654-CAL4

S703654-CAL5

S703654-CAL6

S703654-CAL7

S703654-CAL8

S703654-CAL9

S703654-CALA

S703654-CALB

S703654-ICV1

S703654-LCV1

S703654-LCV2

S703654-TUN1

S704360

Volatile Organic Compounds

S704360-CAL1 S704360-CAL2 S704360-CAL3 S704360-CAL4 S704360-CAL5 S704360-CAL6 S704360-CAL7 S704360-CAL8 S704360-CAL9 S704360-CALA S704360-CALB S704360-ICV1 S704360-LCV1 S704360-LCV2 S704360-LCV3 S704360-LCV4

S704555

Volatile Organic Compounds

S704555-CCV1 S704555-TUN1

S704360-TUN1

S704678

Semivolatile Organic Compounds by GCMS

S704678-CCV1 S704678-TUN1

S704750

Semivolatile Organic Compounds by GCMS

S704750-CCV1 S704750-TUN1